

Brian D Weitzner

List of Publications by Year in descending order

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Version: 2024-02-01

19
papers

2,403
citations

623734

14
h-index

794594

19
g-index

22
all docs

22
docs citations

22
times ranked

3800
citing authors

#	ARTICLE	IF	CITATIONS
1	Macromolecular modeling and design in Rosetta: recent methods and frameworks. <i>Nature Methods</i> , 2020, 17, 665-680.	19.0	513
2	De novo design of potent and selective mimics of IL-2 and IL-15. <i>Nature</i> , 2019, 565, 186-191.	27.8	362
3	Serverification of Molecular Modeling Applications: The Rosetta Online Server That Includes Everyone (ROSIE). <i>PLoS ONE</i> , 2013, 8, e63906.	2.5	348
4	Benchmarking and Analysis of Protein Docking Performance in Rosetta v3.2. <i>PLoS ONE</i> , 2011, 6, e22477.	2.5	272
5	Modeling and docking of antibody structures with Rosetta. <i>Nature Protocols</i> , 2017, 12, 401-416.	12.0	236
6	An Integrated Framework Advancing Membrane Protein Modeling and Design. <i>PLoS Computational Biology</i> , 2015, 11, e1004398.	3.2	145
7	RosettaAntibodyDesign (RABD): A general framework for computational antibody design. <i>PLoS Computational Biology</i> , 2018, 14, e1006112.	3.2	115
8	Blind prediction performance of RosettaAntibody 3.0: Grafting, relaxation, kinematic loop modeling, and full CDR optimization. <i>Proteins: Structure, Function and Bioinformatics</i> , 2014, 82, 1611-1623.	2.6	91
9	The Origin of CDR H3 Structural Diversity. <i>Structure</i> , 2015, 23, 302-311.	3.3	78
10	Real-Time PyMOL Visualization for Rosetta and PyRosetta. <i>PLoS ONE</i> , 2011, 6, e21931.	2.5	55
11	Accurate Structure Prediction of CDR H3 Loops Enabled by a Novel Structure-Based C-Terminal Constraint. <i>Journal of Immunology</i> , 2017, 198, 505-515.	0.8	43
12	Anchor extension: a structure-guided approach to design cyclic peptides targeting enzyme active sites. <i>Nature Communications</i> , 2021, 12, 3384.	12.8	37
13	A computational method for design of connected catalytic networks in proteins. <i>Protein Science</i> , 2019, 28, 2036-2041.	7.6	28
14	Better together: Elements of successful scientific software development in a distributed collaborative community. <i>PLoS Computational Biology</i> , 2020, 16, e1007507.	3.2	27
15	Integration of the Rosetta suite with the python software stack via reproducible packaging and core programming interfaces for distributed simulation. <i>Protein Science</i> , 2020, 29, 43-51.	7.6	13
16	The influence of proline isomerization on potency and stability of anti-HIV antibody 10E8. <i>Scientific Reports</i> , 2020, 10, 14313.	3.3	12
17	A Framework to Simplify Combined Sampling Strategies in Rosetta. <i>PLoS ONE</i> , 2015, 10, e0138220.	2.5	10
18	PyRosetta Jupyter Notebooks Teach Biomolecular Structure Prediction and Design. <i>The Biophysicist</i> , 2021, 2, 108-122.	0.3	8

#	ARTICLE	IF	CITATIONS
19	I am a United Academic Worker. Science, 2017, 358, 266-266.	12.6	1